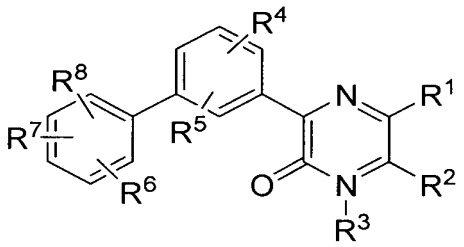
AMENDMENTS TO THE CLAIMS:

This listing of claims replaces all prior versions, and listings, of claims in the application:

LISTING OF CLAIMS:

1. (Original) A compound represented by Formula (I):



(I)

or pharmaceutically acceptable salts thereof, wherein

R^1 and R^2 each independently is
(a) H,
(b) C_1-C_6-alkyl, optionally substituted with one or more substituents selected from the group consisting of: F, CF_3, OH, NR^aR^b, COOH, $CONR^aR^b$, $SO_2NR^aR^b$, $C(=NH)NH_2$, tetrazolyl, triazolyl, oxazolyl, oxadiazolyl, isooxazolyl, thiazolyl, pyrazolyl, pyridyl, pyrimidinyl, pyrazinyl, phenyl, piperidinyl, morpholinyl, pyrrolidinyl and piperazinyl,
(c) $-C(=O)R^a$, $COOR^a$, $CONR^aR^b$,
(d) $-C_0$-C_4-alkyl-C_1-C_4-perfluoroalkyl,
(e) NR^aR^b, $-N(COR^a)R^b$, $-N(SO_2R^a)R^b$, or
(f) tetrazolyl, triazolyl, oxazolyl, oxadiazolyl, isooxazolyl, thiazolyl, pyrazolyl, pyridyl, pyrimidinyl, pyrazinyl, phenyl, piperidinyl, morpholinyl, pyrrolidinyl or piperazinyl, any of which is optionally substituted with 1-3 substituents independently selected from the group consisting of: F, Cl, Br, I and CN;

R^a is
(a) H,

(b) C_1-C_6-alkyl, optionally substituted with one or more substituents independently selected from the group consisting of CF_3 and O-(C_1-C_4)alkyl,
(c) C_0-C_4-alkyl-(C_1-C_4)-perfluoroalkyl,
(d) NH_2,
(e) C_1-C_4-alkyl-phenyl, C_1-C_4-alkyl-pyridyl, or
(f) C_3-C_7-cycloalkyl, optionally substituted with one or more substituents selected from the group consisting of F, Cl, Br, OH, -O-C_1-C_4-alkyl, and C_1-C_4-alkyl;

R^b is
(a) H, or
(b) C_1-C_6-alkyl;

R^3 is:
(a) H,
(b) -C_1-C_4-alkyl, optionally substituted with one or more substituents independently selected from the group consisting of: F, CF_3, Cl, N, OH, O-(C_1-C_4)alkyl, $S(O)_{0-2}$-(C_1-C_4)alkyl, O-$CONR^aR^b$, NR^aR^b, $N(R^a)CONR^aR^b$, $COOR^a$, CN, $CONR^aR^b$, $SO_2NR^aR^b$, $N(R^a)SO_2NR^aR^b$, -C(=NH)NH_2, tetrazolyl, triazolyl, imidazolyl, oxazolyl, oxadiazolyl, isooxazolyl, thiazolyl, furyl, thienyl, pyrazolyl, pyrrolyl, pyridyl, pyrimidinyl, pyrazinyl, phenyl, piperidinyl, morpholinyl, pyrrolidinyl and piperazinyl,
or R^a and R^b, together with N to which they are attached, may form a C_3-C_7-cycloalkyl or a C_3-C_7-heterocycloalkyl, wherein said cycloalkyl and heterocycloalkyl is optionally substituted with one or more substituents selected from the group consisting of: F, Cl, Br, OH, -O-C_1-C_4-alkyl, and C_1-C_4-alkyl,

(c) -C_0-C_4-alkyl-C_1-C_4-perfluoroalkyl, or
(d) C_1-C_4-alkyl-C(=O)-R^a, -C_1-C_4-alkyl-C(=O)-C_1-C_4-perfluoroalkyl, or
(e) -C_1-C_4-alkyl-C_3-C_7-cycloalkyl, wherein said cycloalkyl is optionally substituted with one or more substituents selected from the group consisting of: F, Cl, Br, OH, -O-C_1-C_4-alkyl, and C_1-C_4-alkyl;

R^4 and R^5 each independently is:
(a) H,
(b) -C_1-C_6-alkyl, optionally substituted with one or more substituents independently selected from the group consisting of: F, CF_3 and -O-(C_1-C_4)alkyl,

(c) -O-C_0-C_6-alkyl, -O-phenyl, -O-C_1-C_4-alkyl-phenyl, -O-pyridyl, -O-C_1-C_4-alkyl-pyridyl, wherein phenyl and pyridyl are optionally substituted with 1-3 substituents independently selected from the group consisting of: F, Cl, Br, I and CN,

(d) -C_0-C_4-alkyl-C_1-C_4-perfluoroalkyl, -O-C_0-C_4-alkyl-C_1-C_4-perfluoroalkyl, or

(e) F, Cl, Br, I; and

R^6, R^7 and R^8 each independently is:

(a) H,

(b) C_1-C_6-alkyl,

(c) -O- C_1-C_6-alkyl, optionally substituted with one or more substituents independently selected from the group consisting of: F and CF_3,

(d) -C_0-C_4-alkyl-C_1-C_4-perfluoroalkyl, -O-C_0-C_4-alkyl-C_1-C_4-perfluoroalkyl,

(e) -O-phenyl, -O-C_1-C_4-alkyl-phenyl, -O-pyridyl, -O-C_1-C_4-alkyl-pyridyl, wherein phenyl and pyridyl are optionally substituted with 1-3 substituents independently selected from the group consisting of: F, Cl, Br, I, and CN, or

(f) F, Cl, Br, I, -OR^a, phenyl or pyridyl, wherein phenyl and pyridyl are optionally substituted with one or more substituents independently selected from the group consisting of: F, Cl, Br, I and CN,

with the proviso that when R^6 and R^7 are present on adjacent carbon atoms, R^6 and R^7, together with the benzene ring to which they are attached, may form a bicyclic aromatic ring selected from the group consisting of: naphthyl, quinolinyl and benzothiazolyl, any aromatic ring of which is optionally substituted with 1-4 substituents independently selected from F, Cl, Br, I and CN.

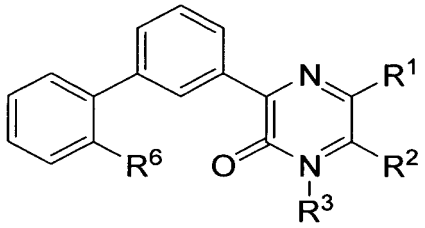
2. (Original) The compound of Claim 1 described by the chemical Formula (I), or a pharmaceutically acceptable salt thereof, wherein

R_6 is other than H and is attached at the ortho position, and all other variables are as previously defined.

3. (Original) The compound of Claim 2, or a pharmaceutically acceptable salt thereof, wherein

R^1 is H, $COOR^a$ or $CONR^aR^b$, and all other variables are as previously defined.

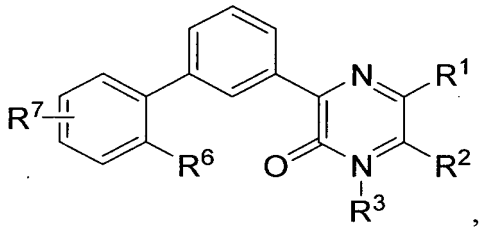
4. (Original) The compound of Claim 1 represented by Formula Ia:



or a pharmaceutically acceptable salt thereof, wherein

R^6 is OR^a or C_0-C_4-alkyl-C_1-C_4-perfluoroalkyl, and all other variables are as previously defined.

5. (Original) The compound of Claim 1 represented by Formula Ib:

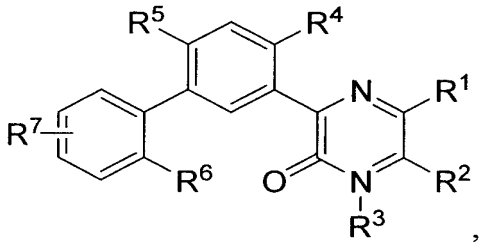


or a pharmaceutically acceptable salt thereof, wherein

R^6 is OR^a or C_0-C_4-alkyl-C_1-C_4-perfluoroalkyl;

R^7 is H, F, Cl, Br or I; and all other variables are as previously defined.

6. (Original) The compound of Claim 1 represented by Formula Ic:



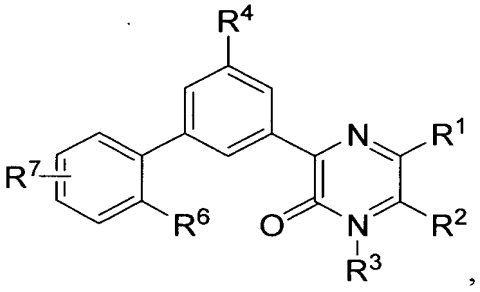
or a pharmaceutically acceptable salt thereof, wherein

R^4 and R^5 each independently is H, F, Cl, Br or I;

R^6 is OR^a or C_0-C_4-alkyl-C_1-C_4-perfluoroalkyl;

R^7 is H, F, Cl, Br or I; and all other variables are as previously defined.

7. (Original) The compound of Claim 1 represented by Formula Id:



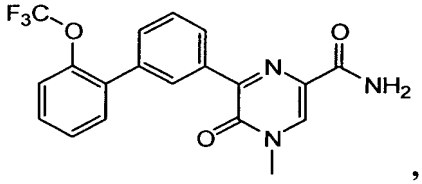
or a pharmaceutically acceptable salt thereof, wherein

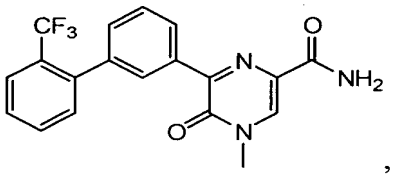
R^4 is F, Cl, Br or I;

R^6 is OR^a or C_0-C_4-alkyl-C_1-C_4-perfluoroalkyl;

R^7 is H, F, Cl, Br or I; and all other variables are as previously defined.

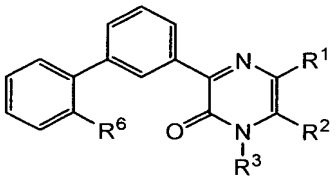
8. (Original) A compound selected from





and pharmaceutically acceptable salts thereof.

9. (Original) A compound according to Claim 1 represented by

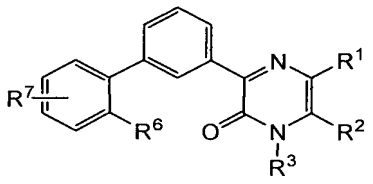


R^6	R^3	R^2	R^1
$-OCF_3$	H	H	$-CONH_2$

R^6	R^3	R^2	R^1
$-OCF_3$	CH_3	H	-COOH
$-OCF_3$	CH_3	$-CONH_2$	H
$-CF_3$	CH_3	$-CONH_2$	H
$-CF_3$	H	H	-COO-t-Bu
$-CF_3$	H	H	-CONH-t-Bu
$-CF_3$	H	H	-COOH
$-OCH_2CF_2CF_3$	H	H	$-CONH_2$
$-OCH_2CF_2CF_3$	H	H	-COOH
$-OCH_2CF_3$	H	H	$-CONH_2$
$-OCH_2CF_3$	CH_3	H	$-CONH_2$
$-OCH_2CF_3$	H	$-CONH_2$	H
$-OCH_2CF_3$	CH_3	$-CONH_2$	H
$-OCH_2CF_2CF_3$	CH_3	H	$-CONH_2$
$-OCH_2CF_3$	H	H	-COOH
$-CF_3$	$-CH_2CF_3$	H	$-CONH_2$
$-CF_3$	$-C(CH_3)_3$	H	$-CONH_2$
$-CF_3$	$-CH(CH_3)_2$	H	$-CONH_2$
$-CF_3$	$-CH_2CH_3$	H	$-CONH_2$
$-OCF_3$	CH_3	$-CONH_2$	$-CONH_2$

or a pharmaceutically acceptable salt thereof.

10. (Original) A compound according to Claim 1 represented by

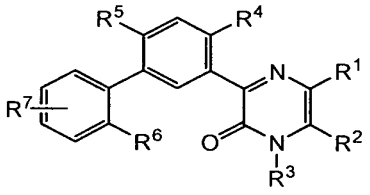


R^7	R^6	R^3	R^2	R^1
$4\text{-}CF_3$	$-CF_3$	$-CH_3$	H	$-CONH_2$
$4\text{-}CF_3$	$-CF_3$	$-CH_3$	$-CONH_2$	H

R^7	R^6	R^3	R^2	R^1
5-F	$-OCF_3$	$-CH_3$	H	$-CONH_2$
$5-CF_3$	$-OCF_3$	$-CH_3$	H	$-CONH_2$
5-F	$-OCF_3$	H	H	$-CONH_2$
5-F	$-OCH_2CF_2CF_3$	H	H	$-CONH_2$
5-F	$-OCH_2CF_3$	H	H	$-CONH_2$
5-F	$-OCH_2CF_2CF_3$	$-CH_3$	H	$-CONH_2$
5-F	$-OCH_2CF_3$	$-CH_3$	H	$-CONH_2$
6-F	$-CF_3$	H	H	$-CONH_2$
3-F	$-CF_3$	H	H	$-CONH_2$
6-F	$-CF_3$	$-CH_3$	H	$-CONH_2$
3-F	$-CF_3$	$-CH_3$	H	$-CONH_2$
5-F	$-CF_3$	$-CH_3$	H	$-CONH_2$
4-F	$-CF_3$	$-CH_3$	H	$-CONH_2$
4-F	$-CF_3$	H	H	$-CONH_2$
5-F	$-CF_3$	H	H	$-CONH_2$
$5-CH_3$	$-CF_3$	$-CH_3$	H	$-CONH_2$
$4-CH_3$	$-CF_3$	$-CH_3$	H	$-CONH_2$

or a pharmaceutically acceptable salt thereof.

11. (Original) A compound according to Claim 1 represented by



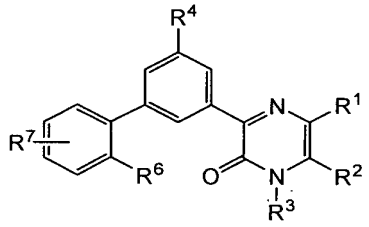
R^7	R^6	R^5	R^4	R^3	R^2	R^1
H	$-OCF_3$	F	H	$-CH_3$	$-CONH_2$	H
H	$-OCF_3$	F	H	$-CH_3$	H	$-CONH_2$
H	$-OCF_3$	H	F	$-CH_3$	H	H

R^7	R^6	R^5	R^4	R^3	R^2	R^1
H	$-OCF_3$	H	F	$-CH_3$	H	$-CONH_2$
H	$-OCH_2CF_3$	H	F	$-CH_3$	H	$-CONH_2$
H	$-OCF_3$	H	F	$-CH_3$	H	-COOH
5-F	$-OCF_3$	F	H	$-CH_3$	H	$-CONH_2$
5-F	$-OCF_3$	H	F	$-CH_3$	H	$-CONH_2$
H	$-CF_3$	H	F	$-CH_3$	H	$-CONH_2$
H	$-CF_3$	F	H	$-CH_3$	H	$-CONH_2$
H	$-OCF_3$	H	F	H	H	$-CONH_2$
H	$-OCF_3$	F	H	H	H	$-CONH_2$
5-F	$-OCF_3$	F	H	H	H	$-CONH_2$
H	$-CF_3$	F	H	$-CH_3$	H	-COOH
5-F	$-OCH_2CF_2CF_3$	F	H	$-CH_3$	H	$-CONH_2$
H	$-CF_3$	H	F	$-CH_3$	H	-COOH
H	$-OCF_3$	F	H	$-CH_3$	H	$-COOCH_3$
H	$-CF_3$	F	H	$-CH_3$	H	$-COOCH_3$
H	$-CF_3$	F	H	$-CH_3$	H	$-COOCH_3$
5-F	$-OCH_2CH_3$	F	H	$-CH_3$	H	$-COOCH_3$
5-F	$-OCH_2CF_3$	F	H	$-CH_3$	H	$-COOCH_3$
H	$-OCF_3$	F	H	H	H	-COOH
H	$-OCF_3$	H	F	H	H	-COOH
5-F	$-OCH_2CF_2CF_3$	F	H	$-CH_3$	H	-COOH
H	$-CF_3$	H	F	H	H	$-CONH_2$
H	$-CF_3$	Br	H	$-CH_3$	H	$-COOCH_3$
H	$-CF_3$	F	H	H	H	$-CONH_2$
H	$-CF_3$	Br	H	$-CH_3$	H	$-CONH_2$
5-F	$-CF_3$	H	F	$-CH_3$	H	$-CONH_2$
H	$-CF_3$	Br	H	H	H	$-COOCH_3$
H	$-CF_3$	Br	H	H	H	-COOH
5-F	$-OCH_2CF_3$	F	H	H	H	$-CONH_2$
3-F	$-CF_3$	F	H	$-CH_3$	H	$-CONH_2$
3-F	$-CF_3$	H	F	$-CH_3$	H	$-CONH_2$
5-F	$-CF_3$	F	H	$-CH_3$	H	$-CONH_2$

R^7	R^6	R^5	R^4	R^3	R^2	R^1
4-F	$-CF_3$	F	H	$-CH_3$	H	$-CONH_2$
4-F	$-CF_3$	H	F	H	H	$-CONH_2$

or a pharmaceutically acceptable salt thereof.

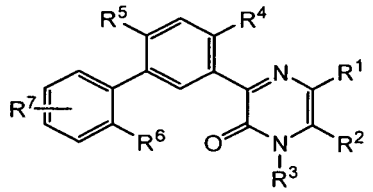
12. (Original) A compound according to Claim 1 represented by

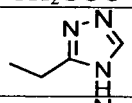
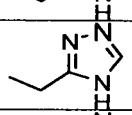
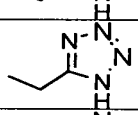
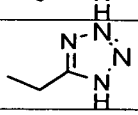


R^7	R^6	R^4	R^3	R^2	R^1
H	$-OCF_3$	F	$-CH_3$	H	$-CONH_2$
H	$-OCF_3$	F	$-CH_3$	$-CONH_2$	H
H	$-OCF_3$	F	$-CH_3$	H	-COOH
H	$-OCF_3$	F	H	H	$-CONH_2$
H	$-OCF_3$	F	H	H	-COOH
H	$-CF_3$	F	$-CH_3$	H	$-CONH_2$
H	$-CF_3$	F	$-CH_3$	H	-COOH
5-F	$-OCH_2CF_2CF_3$	F	$-CH_3$	H	$-CONH_2$
5-F	$-OCH_2CF_3$	F	$-CH_3$	H	$-CONH_2$
H	$-OCF_3$	F	$-CH_3$	H	-COOCH3
5-F	$-CF_3$	F	$-CH_3$	H	$-CONH_2$
H	$-CF_3$	F	H	H	$-CONH_2$
3-F	$-CF_3$	F	$-CH_3$	H	$-CONH_2$

or a pharmaceutically acceptable salt thereof.

13. (Original) A compound according to Claim 1 represented by

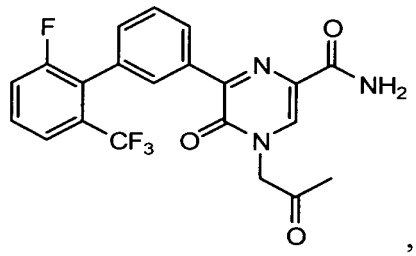


R^7	R^6	R^5	R^4	R^3	R^2	R^1
H	OCF_3	H	H	CH_2CONH_2	H	$CONH_2$
H	CF_3	H	H	CH_2CONH_2	H	$CONH_2$
H	OCF_3	H	H	CH_2COOH	H	$CONH_2$
H	OCF_3	H	H	CH_2COO-tBu	H	$CONH_2$
H	OCF_3	H	H	CH_2CN	H	$CONH_2$
H	CF_3	H	H	CH_2CN	H	$CONH_2$
H	CF_3	H	H	CH_2COOH	H	$CONH_2$
H	CF_3	H	H	CH_2COO-tBu	H	$CONH_2$
H	OCF_3	H	H		H	$CONH_2$
H	CF_3	H	H		H	$CONH_2$
H	CF_3	H	H		H	$CONH_2$
H	OCF_3	H	H		H	$CONH_2$
H	OCF_3	H	H	CH_2CH_2OH	H	$CONH_2$
H	CF_3	H	H	CH_2CH_2OH	H	$CONH_2$
H	OCF_3	H	H	$CH_2CH_2NH_2$	H	$CONH_2$
H	CF_3	H	H	$CH_2CH_2NH_2$	H	$CONH_2$
H	OCF_3	H	H	$CH_2CH_2N(CH_3)_2$	H	$CONH_2$
H	CF_3	H	H	$CH_2CH_2N(CH_3)_2$	H	$CONH_2$
4-F	CF_3	H	H	CH_2CONH_2	H	$CONH_2$
5-F	CF_3	H	H	CH_2CONH_2	H	$CONH_2$
H	CF_3	F	H	CH_2CONH_2	H	$CONH_2$
H	CF_3	F	H	CH_2CONH_2	H	$CONH_2$
R^7	R^6	R^5	R^4	R^3	R^2	R^1

4-F	CF_3	F	H	CH_2CONH_2	H	$CONH_2$
5-F	CF_3	F	H	CH_2CONH_2	H	$CONH_2$

or a pharmaceutically acceptable salt thereof.

14. (Original) A compound selected from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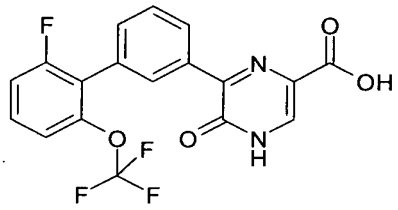
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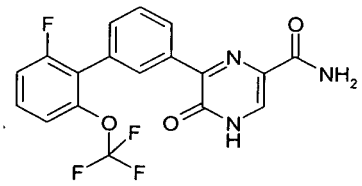
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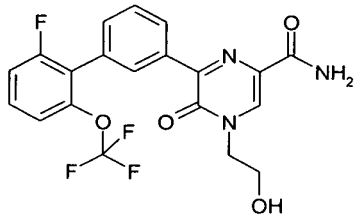
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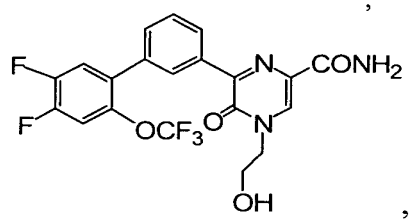
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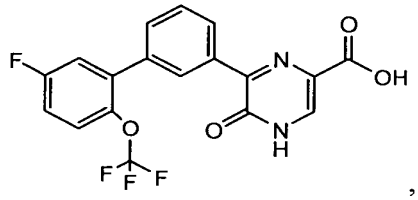
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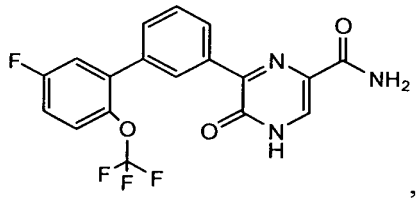
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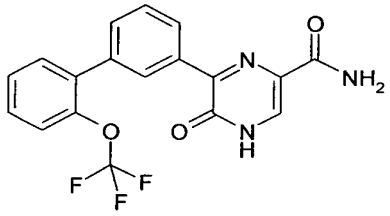
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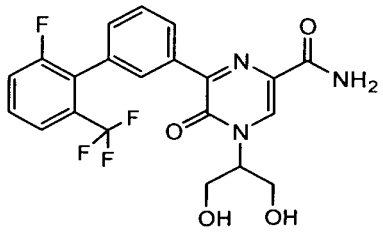
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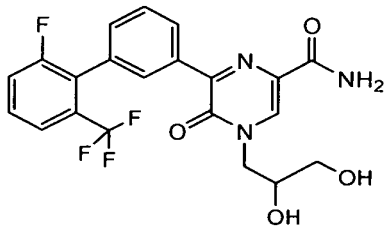
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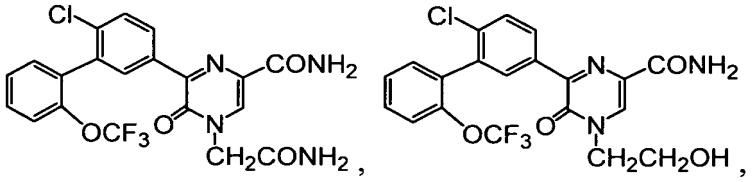
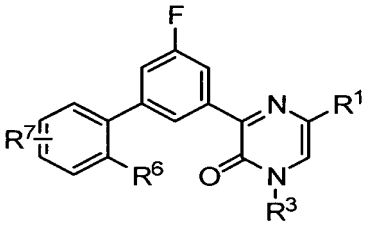


and pharmaceutically acceptable salts thereof.

15. (Original) A compound according to Claim 1 represented by

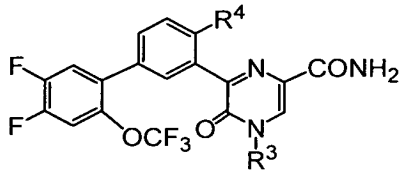


R^7	R^6	R^3	R^1
4-F	$-CF_3$	$-CH_3$	$-CONH_2$
5-F	$-CF_3$	H	$-CONH_2$
H	CF_3	$-CH_2CH_2OH$	$-CONH_2$
H	$-CF_3$	$-CH_2CH_2F$	$-CONH_2$
H	$-OCF_3$	$-CH_2CH_2OH$	$-CONH_2$
H	$-OCF_3$	$-CH_2CH_2CH_2OH$	$-CONH_2$
H	$-OCF_3$	$-CH_2CONH_2$	$-CONH_2$
H	$-OCF_3$	$-CH(CH_3)CONH_2$	$-CONH_2$
H	$-OCF_3$	$-CH_2CH_3$	$-CONH_2$
H	$-OCF_3$	$-CH_2CH_2F$	$-CONH_2$
H	$-CF_3$	$-CH_2CONH_2$	$-CONH_2$
H	$-OCF_3$	$-CH_2SO_2NH_2$	$-CONH_2$
3-F	$-CF_3$	$-CH_2SO_2NH_2$	$-CONH_2$
H	$-CF_3$	$-CH_2CH_3$	$-CONH_2$
H	$-CF_3$	$-CH_2CH_2CH_2OH$	$-CONH_2$
H	$-OCF_3$	$-CH_2CH_2Cl$	$-CONH_2$
H	$-OCF_3$	$-CH_2CH_2N_3$	$-CONH_2$
H	$-OCF_3$	$-CH_2CH_2NH_2$	$-CONH_2$

R^7	R^6	R^3	R^1
H	$-CF_3$	$-CH(CH_3)CONH_2$	$-CONH_2$
H	$-OCF_3$	$-CH_2CONHCH_3$	$-CONH_2$
4-F	$-CF_3$	$-CH_2CONH_2$	$-CONH_2$
4-F	$-CF_3$	$-CH_2CH_2OH$	$-CONH_2$
4-F	$-CF_3$	$-CH_2CH_3$	$-CONH_2$
4-F	$-CF_3$	H	$-CONH_2$
5-F	$-CF_3$	$-CH_2CH_2OH$	$-CONH_2$
5-F	$-CF_3$	$-CH_2CONH_2$	$-CONH_2$
H	$-OCF_3$	$-CH_2C(=O)CH_3$	$-CONH_2$

or a pharmaceutically acceptable salt thereof.

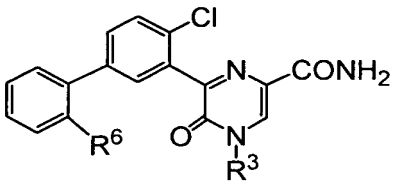
16. (Original) A compound according to Claim 1 represented by



R^4	R^3
F	H
F	CH_2CONH_2
F	CH_2CH_2OH
F	CH_3
H	H
H	CH_2CONH_2
H	CH_3

or a pharmaceutically acceptable salt thereof.

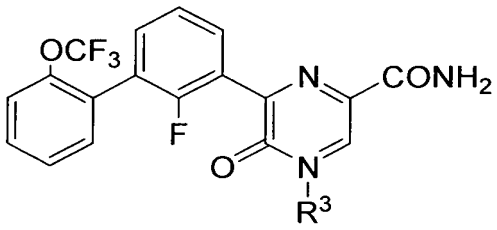
17. (Original) A compound according to Claim 1 represented by



R^6	R^3
OCF3	CH_2CONH_2
OCF3	$CH(CH_3)CONH_2$
OCF3	CH_2CH_2OH
OCF3	$CH_2CH_2CH_2OH$
OCF3	CH_2CH_3
CF3	CH_2CONH_2
CF3	$CH(CH_3)CONH_2$
CF3	CH_2CH_2OH
CF3	$CH_2CH_2CH_2OH$
CF3	CH_2CH_3

or a pharmaceutically acceptable salt thereof.

18. (Original) A compound according to Claim 1 represented by

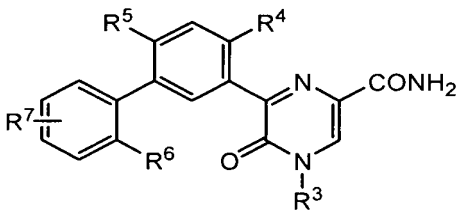


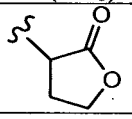
R^3
H
CH_2CONH_2
CH_2CH_2OH
$CH(CH_3)CONH_2$
$CH_2CH_2CH_2OH$
CH_2CH_3
$CH_2CONHCH_3$

R^3
$CH_2CON(CH_3)_2$
$CH_2CO_2CH_3$

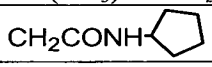
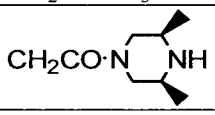
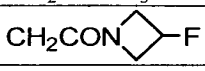
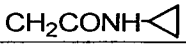
or a pharmaceutically acceptable salt thereof.

19. (Original) A compound according to Claim 1 represented by



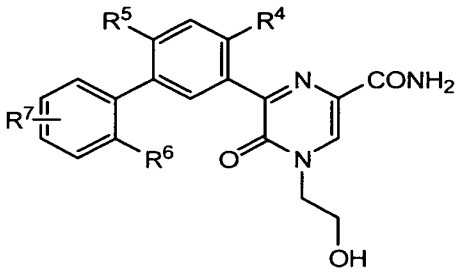
R^7	R^6	R^5	R^4	R^3
H	OCF_3	H	H	$C(CH_3)_2CONH_2$
5-Cl	Cl	H	H	CH_2CONH_2
4- CF_3	CF_3	H	H	CH_2COOH
4- CF_3	CF_3	H	H	CH_2CONH_2
H	Cl	H	H	CH_2CONH_2
3-Cl	Cl	H	H	CH_2CONH_2
5- CF_3	CF_3	H	H	CH_2CONH_2
H	OCF_3	H	H	$CH(CH_3)COOH$
H	OCF_3	H	H	$CH(CH_3)CONH_2$
H	CF_3	H	H	$CH(CH_3)CONH_2$
H	OCF_3	H	H	
R^7	**R^6**	**R^5**	**R^4**	**R^3**
H	OCF_3	H	H	$CH(CH_3)COCH_3$
H	OCF_3	H	F	CH_2CONH_2
6-F	CF_3	H	H	CH_2CONH_2
H	OCF_3	F	H	CH_2CONH_2
6-F	CF_3	H	H	$CH(CH_3)CONH_2$

H	OCF_3	F	H	$CH(CH_3)CONH_2$
4-F	OCF_3	F	H	CH_2CONH_2
4-F	OCF_3	F	H	$CH(CH_3)CONH_2$
H	CF_3	H	F	CH_2CONH_2
3-F	CF_3	H	H	CH_2CONH_2
4-Cl	CF_3	H	H	CH_2CONH_2
H	OCF_3	H	F	$CH(CH_3)CONH_2$
6-F	CF_3	H	H	$CH_2CONHCH_3$
6-F	CF_3	H	H	$CH_2CON(CH_3)_2$
H	CF_3	H	H	$CH(CONH_2)_2$
H	OCF_3	H	H	$CH(CONH_2)_2$
6-F	CF_3	H	H	$CH(CONH_2)_2$
H	CF_3	H	H	$CH_2CONHCH_3$
H	CF_3	H	H	$CH_2CON(CH_3)_2$
$4\text{-}CF_3$	CF_3	H	F	CH_2CONH_2
H	OCF_3	H	H	$CH_2CON(CH_3)_2$
H	OCF_3	H	H	$CH_2CONHCH_3$
H	OCF_3	F	H	$CH_2CONHCH_3$
H	OCF_3	H	F	$CH_2CONHCH_3$
6-F	CF_3	H	H	$CH_2CONH(CH_2)_2OH$
H	CF_3	H	F	$CH_2CONHCH_3$
H	OCH_2CF_3	H	H	CH_2CONH_2
H	$OCH_2CF_2CF_3$	H	H	CH_2CONH_2
6-F	CF_3	H	H	CH_2CONH 
H	CF_3	H	H	CH_2CONH 
6-F	CF_3	H	H	$CH_2CONHCH(CH_3)_2$
6-F	CF_3	H	H	$CH_2CONHC(CH_3)_3$
R^7	**R^6**	**R^5**	**R^4**	**R^3**
H	CF_3	H	H	$CH_2CONHC(CH_3)_3$
4-F	OCF_3	H	H	CH_2CONH_2
H	CF_3	H	H	$CH_2CONHCH_2CH_3$
H	CF_3	H	F	$CH_2CONHCH_2CH_3$
6-F	CF_3	H	H	CH_2CON 

6-F	OCF_3	H	H	CH_2CONH_2
6-F	CF_3	H	H	
5-F	OCF_3	H	H	CH_2CONH_2
6-F	OCF_3	H	H	CH_2COOH
H	OCF_3	H	OCH_2Ph	CH_2CONH_2
H	OCF_3	H	OCH_2Ph	$CH(CH_3)CONH_2$
H	CF_3	H	H	
5-I	OCF_3	H	H	CH_2CONH_2
6-F	CF_3	H	H	CH_2COCH_3
H	CF_3	H	H	CH_2COCH_3
H	CF_3	H	H	
H	CF_3	I	I	CH_2CONH_2
5-F	OCF_3	H	F	CH_2CONH_2
4-F	CF_3	H	F	CH_2CONH_2
H	OCF_3	H	H	CH_2COCH_3
H	CF_3	H	H	
H	OCF_3	Br	H	CH_2CONH_2
H	OCF_3	H	F	CH_2COCH_3
5-F	CF_3	H	H	CH_2COCH_3
5-F	CF_3	H	H	
H	OCF_3	F	H	CH_2COCH_3

or a pharmaceutically acceptable salt thereof.

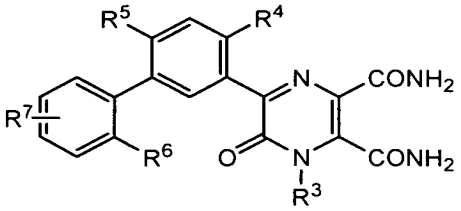
20. (Original) A compound according to Claim 1 represented by



R^7	R^6	R^5	R^4
4-F	CF_3	H	H
6-F	CF_3	H	H
3-F	CF_3	H	H
H	CF_3	F	H
H	CF_3	H	F
5-CF_3	OCF_3	H	H
4- F	OCF_3	H	H
H	OCF_3	F	H
H	OCF_3	H	F
4-F	CF_3	H	F
3-F	CF_3	H	F
H	O $CH_2CF_2CF_3$	F	H
4-F	OCF_3	F	H
4-CF_3	CF_3	H	H
5-Cl	Cl	H	H
5-CF_3	CF_3	H	H
5-F	CF_3	H	H
4-F	OCF_3	H	H
5-F	OCF_3	H	H
H	OCF_3	H	OCH_2Ph
H	OCF_3	Br	H
4-Cl	CF_3	H	H

or a pharmaceutically acceptable salt thereof.

21. (Original) A compound according to Claim 1 represented by

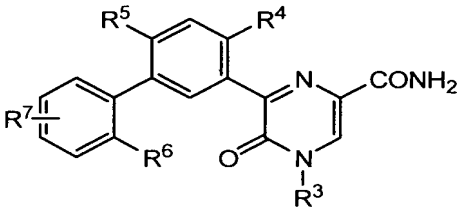


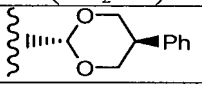
R^7	R^6	R^5	R^4	R^3
H	OCF_3	H	H	CH_3
H	CF_3	H	H	CH_3
H	CF_3	H	H	H
H	OCF_3	H	H	H
4-F	CF_3	H	H	H
4-F	CF_3	H	H	CH_3
4-F	CF_3	H	H	CH_2CH_2OH
3-F	CF_3	H	H	H
3-F	CF_3	H	H	CH_3
3-F	CF_3	H	H	CH_2CH_2OH
4-Cl	CF_3	H	H	CH_3
5- CF_3	CF_3	H	H	CH_3
4-Cl	CF_3	H	H	H
5- CF_3	CF_3	H	H	H
4-F	OCF_3	H	H	H
4-F	OCF_3	H	H	CH_3
4-F	OCF_3	H	H	CH_2CH_2OH
4-Cl	CF_3	H	H	CH_2CH_2OH
5- CF_3	CF_3	H	H	CH_2CH_2OH
H	$OCH_2CF_2CF_3$	H	H	H
H	$OCH_2CF_2CF_3$	H	H	CH_3
H	$OCH_2CF_2CF_3$	H	H	CH_2CH_2OH
H	OCH_2CF_3	H	H	H
H	OCF_3	F	H	H
H	CF_3	F	H	H
R^7	**R^6**	**R^5**	**R^4**	**R^3**
H	OCH_2CF_3	H	H	CH_3

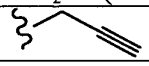
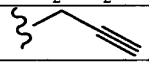
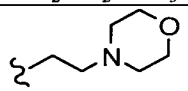
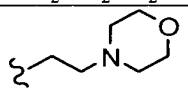
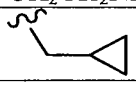
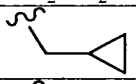
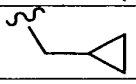
H	OCF_3	F	H	CH_3
H	CF_3	F	H	CH_3
H	OCF_3	F	H	CH_2CH_2OH
H	CF_3	F	H	$CH_2CH{=}CH_2$
H	OCF_3	F	H	$CH_2CH{=}CH_2$
H	CF_3	F	H	$CH_2CH{=}CH_2$
H	CF_3	F	H	$CH_2CH_2OCH_3$

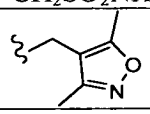
or a pharmaceutically acceptable salt thereof.

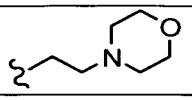
22. (Original) A compound according to Claim 1 represented by



R^7	R^6	R^5	R^4	R^3
H	OCF_3	H	H	$CH_2CH_2OCH_3$
6-F	CF_3	H	H	$CH_2CH_2OCH_3$
6-F	CF_3	H	H	$CH_2CH_2OCH_2CH_3$
H	OCF_3	H	H	$CH_2CH_2CH_3$
H	OCF_3	H	H	$CH(CH_2OH)CH_2OH$
H	CF_3	H	H	$CH(CH_2OH)CH_2OH$
H	OCF_3	H	H	
6-F	OCF_3	H	H	$CH(CH_2OH)CH_2OH$
H	CF_3	H	F	$CH(CH_2OH)CH_2OH$
H	CF_3	H	F	$CH_2CH(OH)CH_2OH$
H	OCF_3	H	H	$CH_2CH_2CH_2CH_3$
H	CF_3	H	H	$CH_2CH(OH)CH_3$
R^7	**R^6**	**R^5**	**R^4**	**R^3**
6-F	CF_3	H	H	$CH_2CH(OH)CH_3$

6-F	OCF_3	H	H	$CH_2CH(OH)CH_2OH$
H	OCF_3	H	H	
H	OCF_3	H	H	$CH_2CH_2CH_2CH_2CH_3$
6-F	CF_3	H	H	
H	OCF_3	H	H	$CH_2C(CH_3)_2CH_2OH$
H	OCF_3	H	H	$CH_2CH_2CONH_2$
H	OCF_3	H	H	$CH_2CH_2SCH_3$
6-F	CF_3	H	H	$CH_2CH_2SCH_3$
H	OCF_3	H	H	
H	OCF_3	H	H	$CH_2CH_2SO_2CH_3$
6-F	CF_3	H	H	$CH_2CH_2SO_2CH_3$
6-F	CF_3	H	H	
6-F	CF_3	H	H	$CH_2CH_2CH_2OH$
6-F	CF_3	H	H	CH_2SCH_3
6-F	CF_3	H	H	$CH_2CH_2CH_2SCH_3$
6-F	CF_3	H	H	$CH_2CH_2CH_2SO_2CH_3$
H	OCF_3	H	H	$CH_2CH_2NHCONH_2$
6-F	CF_3	H	H	
H	OCF_3	H	H	CH_2CF_3
H	OCF_3	H	H	$CH_2CF_2CF_3$
H	CF_3	H	H	$CH_2CH_2CH_2OH$
H	CF_3	H	H	
H	OCF_3	H	H	
H	CF_3	H	H	CH_2SCH_3
H	OCF_3	H	H	CH_2SCH_3
H	OCF_3	H	H	$CH_2SO_2CH_3$
6-F	CF_3	H	H	$CH_2SO_2CH_3$
R^7	**R^6**	**R^5**	**R^4**	**R^3**
6-F	CF_3	H	H	CH_2CH_3

6-F	CF_3	H	H	$CH_2CH_2CH_3$
H	OCF_3	H	H	$CH_2SO_2NHC(CH_3)_3$
H	OCF_3	H	H	$CH_2SO_2NH_2$
H	OCF_3	H	H	
4-F	CF_3	H	H	$CH_2SO_2NH_2$
H	CF_3	H	H	$CH_2SO_2NH_2$
4-F	OCF_3	H	H	$CH_2SO_2NH_2$
6-F	CF_3	H	H	$CH_2SO_2NH_2$
H	CF_3	H	H	$CH_2SO_2CH_3$
H	OCF_3	F	H	$CH_2SO_2CH_3$
4-F	CF_3	H	H	CH_2CH_3
4-F	OCF_3	H	H	$CH_2CH_2CH_2OH$
4-Cl	CF_3	H	H	$CH_2SO_2NH_2$
3-F	CF_3	H	H	$CH_2SO_2NH_2$
4-F	CF_3	H	H	$CH_2SO_2NHCH_3$
H	CF_3	H	F	$CH_2CH_2CH_2OH$
H	OCF_3	F	H	CH_2CH_2F
H	OCF_3	H	F	CH_2CH_2F
4-F	OCF_3	F	H	CH_2CH_2F
6-F	CF_3	H	H	CH_2CH_2F
H	CF_3	H	F	CH_2CH_2F
4- CF_3	CF_3	H	H	$CH_2SO_2NH_2$
H	OCF_3	H	F	$CH_2SO_2NH_2$
H	CF_3	H	F	$CH_2SO_2NH_2$
H	OCF_3	H	F	$CH_2SO_2NHCH_3$
H	CF_3	H	F	$CH_2SO_2NHCH_3$
H	CF_3	H	H	CH_2CH_2F
H	OCF_3	H	H	CH_2CH_2F
3-F	CF_3	H	H	CH_2CH_2F
4-Cl	CF_3	H	H	CH_2CH_2F
R^7	**R^6**	**R^5**	**R^4**	**R^3**
5-CF_3	CF_3	H	H	CH_2CH_2F

H	OCF_3	F	H	
H	OCF_3	H	F	CH_2CH_2F
H	OCF_3	F	H	CH_2CH_2F
5-F	CF_3	H	H	CH_2CH_3
5-F	CF_3	H	H	CH_2CH_2F
H	CF_3	H	H	$CH(CONH_2)_2$
H	OCF_3	H	H	$CH(CONH_2)_2$
6-F	CF_3	H	H	$CH(CONH_2)_2$
6-F	CF_3	H	H	$CH_2CONHCH_2CH_2OH$

or a pharmaceutically acceptable salt thereof.

23. (Original) A pharmaceutical composition comprising a therapeutically effective amount of a compound according to Claim 1, or a pharmaceutically acceptable salt thereof, and a pharmaceutically acceptable carrier.

24. (Cancelled)

25. (Original) A method of treatment or prevention of pain comprising the step of administering to a patient in need thereof a therapeutically effective amount, or a prophylactically effective amount, of a compound according to Claim 1, or a pharmaceutically acceptable salt thereof.

26 to 37. (Cancelled)

38. (Original) A pharmaceutical composition comprising a therapeutically effective amount of a compound according to Claim 2, or a pharmaceutically acceptable salt thereof, and a pharmaceutically acceptable carrier.

39. (Cancelled)

40. (Original) A method of treatment or prevention of pain comprising the step of administering to a patient in need thereof a therapeutically effective amount, or a prophylactically effective amount, of a compound according to Claim 2, or a pharmaceutically acceptable salt thereof.

41 to 52. (Cancelled)